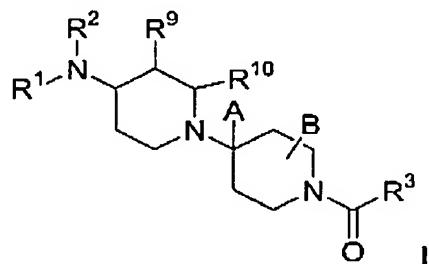


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Presently Pending Claims

Language to be added has been bolded and underlined, while language to be deleted has been ~~bolded and striken-through~~.

21. (Amended) A compound represented by the structural formula I



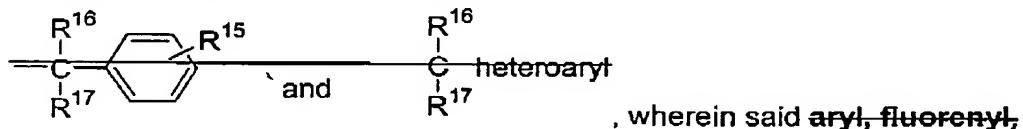
or a pharmaceutically acceptable salt or solvate thereof; wherein:

R<sup>1</sup> is

~~§—M—R<sup>4</sup>~~ ;

R<sup>2</sup> is selected from the group consisting of H, alkyl, aryl, arylalkyl, heteroarylalkyl, alkylketone, arylketone, alkyl, haloalkyl, cycloalkyl, ~~cyclo~~heteroalkyl, cycloalkylalkyl, alkylsulfonyl, arylsulfonyl, alkoxyalkyl, or amide;

R<sup>3</sup> is selected from the group consisting of aryl, 6-membered heteroaryl, fluorenyl; and diphenylmethyl, 6-membered heteroaryl-N-oxide,



R<sup>4</sup> is 1-3 substituents selected from the group consisting of H, halo, alkyl, haloalkyl, alkoxy, cycloalkyl, amide, CF<sub>3</sub>, OCF<sub>3</sub>, aryl, heteroaryl, -XR<sup>7</sup>, -CN, -CO<sub>2</sub>H, -CO<sub>2</sub>R<sup>22</sup>, R<sup>8</sup>-aryl(C<sub>1</sub>-C<sub>6</sub>)alkyl-, R<sup>8</sup>-heteroaryl(C<sub>1</sub>-C<sub>6</sub>)alkyl-, -

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$\text{C}(\text{O})\text{NR}^{21}\text{R}^{22}$ ,  $-\text{C}(\text{O})\text{NH}_2$ , wherein  $\text{R}^4$  can be the same or different and is independently selected when there is more than one  $\text{R}^4$  present;

~~$\text{R}^6$  is selected from the group consisting of H, arylalkyl,  $(\text{C}_1-\text{C}_6)$ alkyl,  $\text{R}^8$ -aryl( $\text{C}_1-\text{C}_6$ )alkyl,  $\text{R}^8$ -heteroaryl( $\text{C}_1-\text{C}_6$ )alkyl,  $-\text{SO}_2-(\text{C}_1-\text{C}_6)$ alkyl,  $-\text{SO}_2-(\text{C}_3-\text{C}_6)$ cycloalkyl,  $-\text{SO}_2$ -aryl,  $\text{R}^8$ -aryl- $\text{SO}_2$ ,  $-\text{C}(\text{O})-(\text{C}_1-\text{C}_6)$ alkyl,  $-\text{C}(\text{O})-(\text{C}_4-\text{C}_6)$ cycloalkyl,  $\text{R}^8$ -aryl- $\text{C}(\text{O})$ ,  $-\text{C}(\text{O})\text{NR}^{21}\text{R}^{22}$ , and  $-\text{SO}_2\text{NR}^{21}\text{R}^{22}$ ;~~

~~$\text{R}^6$  is H,  $(\text{C}_1-\text{C}_6)$ alkyl, or  $(\text{C}_1-\text{C}_6)$ haloalkyl;~~

$\text{R}^7$  is selected from the group consisting of aryl, substituted aryl, heteroaryl, alkyl, haloalkyl and cycloalkyl;

$\text{R}^8$  is 1, 2 or 3 substituents selected from the group consisting of H, halo,  $(\text{C}_1-\text{C}_6)$ alkyl,  $(\text{C}_1-\text{C}_6)$ alkoxy,  $-\text{CF}_3$ ,  $-\text{OCF}_3$ ,  $\text{CH}_3\text{C}(\text{O})$ -,  $-\text{CN}$ ,  $\text{CH}_3\text{SO}_2$ -,  $\text{CF}_3\text{SO}_2$ - and  $-\text{NH}_2$ , wherein  $\text{R}^8$  can be the same or different and is independently selected when there are more than one  $\text{R}^8$  present;

$\text{R}^9$ ,  $\text{R}^{10}$  and B can be the same or different and are each independently selected from the group consisting of hydrogen,  $(\text{C}_1-\text{C}_6)$ alkyl, and  $-(\text{C}_1-\text{C}_6)$ haloalkyl;

$\text{R}^{11}$  and  $\text{R}^{12}$  can be the same or different and are each independently selected from the group consisting of  $(\text{C}_1-\text{C}_6)$ alkyl,  $-(\text{C}_1-\text{C}_6)$ haloalkyl, halogen,  $-\text{NR}^{19}\text{R}^{20}$ ,  $-\text{OH}$ ,  $\text{CF}_3$ ,  $-\text{OCH}_3$ ,  $-\text{O-acyl}$ , and  $-\text{OCF}_3$ ;

$\text{R}^{13}$  is selected from the group consisting of hydrogen,  $\text{R}^{11}$ , H, phenyl,  $-\text{NO}_2$ ,  $-\text{CN}$ ,  $-\text{CH}_2\text{F}$ ,  $-\text{CHF}_2$ ,  $-\text{CHO}$ ,  $-\text{CH}=\text{NOR}^{19}\text{R}_{49}$ , pyridyl-N-oxide, pyrimidinyl, pyrazinyl,  $\text{N}(\text{R}^{20}\text{R}_{29})\text{CONR}^{20}\text{R}^{21}\text{R}_{20}\text{R}_{24}$ ,  $-\text{NHCONH}(\text{chloro-}(\text{C}_1-\text{C}_6)\text{alkyl})$ ,  $-\text{NHCONH}((\text{C}_3-\text{C}_{10})\text{-cycloalkyl}(\text{C}_1-\text{C}_6)\text{alkyl})$ ,  $-\text{NHCO}(\text{C}_1-\text{C}_6)\text{alkyl}$ ,  $-\text{NHCOCF}_3$ ,  $-\text{NHCOCF}_3$ ,  $-\text{NHSO}_2\text{N}((\text{C}_1-\text{C}_6)\text{alkyl})_2$ ,  $-\text{NHSO}_2(\text{C}_1-\text{C}_6)\text{alkyl}$ ,  $-\text{N}(\text{SO}_2\text{CF}_3)_2$ ,  $-\text{NHCO}_2(\text{C}_1-\text{C}_6)\text{alkyl}$ ,  $(\text{C}_3-\text{C}_{10})\text{cycloalkyl}$ ,  $-\text{SR}^{22}$ ,  $-\text{SOR}^{22}$ ,  $-\text{SO}_2\text{R}^{22}$ ,  $-\text{SO}_2\text{NH}(\text{C}_1-\text{C}_6)\text{alkyl}$ ,  $-\text{OSO}_2(\text{C}_1-\text{C}_6)\text{alkyl}$ ,  $-\text{OSO}_2\text{CF}_3$ , hydroxy( $\text{C}_1-\text{C}_6$ )alkyl,  $-\text{CONR}^{19}\text{R}^{20}$ ,  $-\text{CON}(\text{CH}_2\text{CH}_2\text{-O-CH}_3)_2$ ,  $-\text{OCONH}(\text{C}_1-\text{C}_6)\text{alkyl}$ ,  $-\text{CO}_2\text{R}^{19}\text{R}_{49}$ ,  $-\text{Si}(\text{CH}_3)_3$  and  $-\text{B}(\text{OC}(\text{CH}_3)_2)_2$ ;

$\text{R}^{14}$  is selected from the group consisting of  $(\text{C}_1-\text{C}_6)$ alkyl,  $-(\text{C}_1-\text{C}_6)$ haloalkyl,  $-\text{NH}_2$  and  $\text{R}^{15}$ -phenyl;

$\text{R}^{15}$  is 1-3 substituents selected from the group consisting of hydrogen,  $(\text{C}_1-\text{C}_6)$ alkyl,  $-(\text{C}_1-\text{C}_6)$ haloalkyl,  $-\text{CF}_3$ ,  $-\text{CO}_2\text{R}^{20}$ ,  $-\text{CN}$ ,  $(\text{C}_1-\text{C}_6)$ alkoxy and halogen; wherein  $\text{R}^{15}$  can be the same or different and is independently selected when there are more than one  $\text{R}^{15}$  present;

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~~R<sup>16</sup> and R<sup>17</sup> can each be the same or different and are each independently selected from the group consisting of hydrogen and (C<sub>1</sub>-C<sub>6</sub>)alkyl, or~~

~~R<sup>16</sup> and R<sup>17</sup> together are a C<sub>2</sub>-C<sub>5</sub> alkylene group and with the carbon to which they are attached from a spiro ring of 3 to 6 carbon atoms;~~

R<sup>19</sup>, R<sup>20</sup> and R<sup>21</sup> can each be the same or different and are each independently selected from the group consisting of H, (C<sub>1</sub>-C<sub>6</sub>)alkyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl;

R<sup>22</sup> is selected from the group consisting of (C<sub>1</sub>-C<sub>6</sub>)alkyl, -(C<sub>1</sub>-C<sub>6</sub>)haloalkyl, (C<sub>2</sub>-C<sub>6</sub>)hydroxyalkyl, (C<sub>2</sub>-C<sub>6</sub>)alkylene, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl and aryl(C<sub>1</sub>-C<sub>6</sub>)alkyl-;

A is selected from the group consisting of H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, and (C<sub>2</sub>-C<sub>6</sub>)alkenyl.

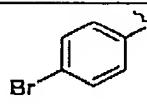
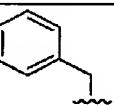
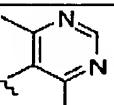
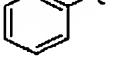
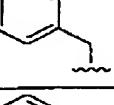
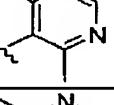
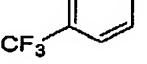
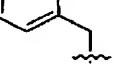
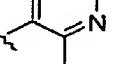
M is aryl or heteroaryl optionally substituted with R<sup>4</sup>;

Q is CH or N; and

X is selected from the group consisting of CH<sub>2</sub>, SO<sub>2</sub>, SO, S, and O, with the following proviso:

when R<sup>1</sup> is phenyl, ~~pyridyl, thiophenyl~~ or naphthyl, R<sup>2</sup> cannot be H, -(C<sub>1</sub>-C<sub>6</sub>)alkyl or -C(O)-(C<sub>1</sub>-C<sub>6</sub>)alkyl.

22. (Amended) A compound having the structural formula I according to claim 21 wherein R<sup>9</sup>, R<sup>10</sup> and B are H, A is -CH<sub>3</sub>, and R<sup>1</sup>, R<sup>2</sup> and R<sup>3</sup> are as defined in the following table:

#	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>
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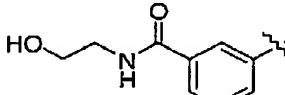
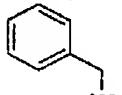
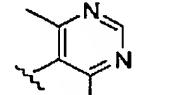
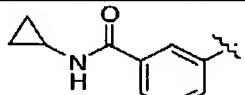
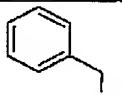
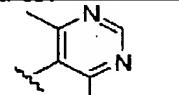
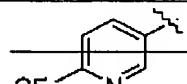
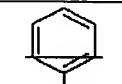
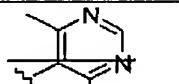
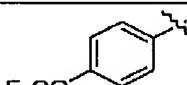
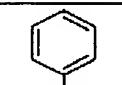
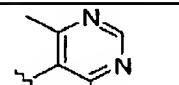
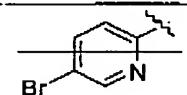
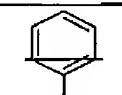
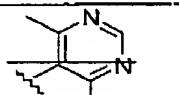
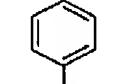
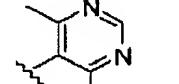
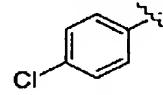
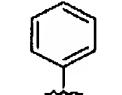
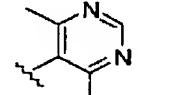
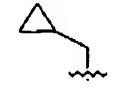
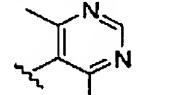
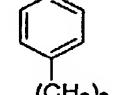
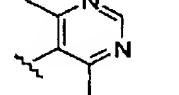
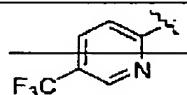
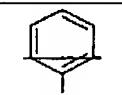
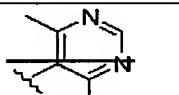
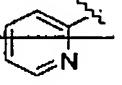
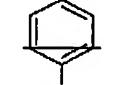
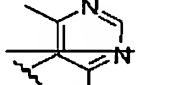
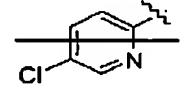
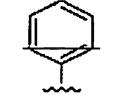
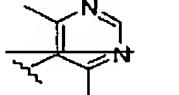
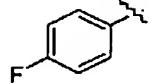
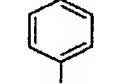
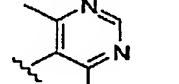
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34	<chem>FC(F)(F)c1ccc(C(F)(F)F)cc1</chem>	<chem>Cc1ccccc1</chem>	<chem>CC1=CC=CC=C1</chem>
35	<chem>Clc1ccc(C(F)(F)F)cc1</chem>	<chem>Cc1ccccc1</chem>	<chem>CC1=CC=CC=C1</chem>
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46	<chem>CC1=CC=CC=C1</chem>	<chem>Cc1ccccc1</chem>	<chem>CC1=CC=CC=C1</chem>

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50	<chem>CC1=CC=C(C=C1)C2=C(C=C(C=C2)C(=O)OC3=CC=CC=C3)C=C3C=CC=C3</chem>	<chem>CC1=CC=C(C=C1)C2=C(C=C(C=C2)C=C3C=CC=C3)C=C3C=CC=C3</chem>	<chem>CC1=CC=C(C=C1)C2=C(C=C(C=C2)C=C3C=CC=C3)C=C3C=CC=C3</chem>
51	<chem>CC1=CC=C(C=C1)C2=C(C=C(C=C2)Cl)C=C3C=CC=C3</chem>	<chem>CC1=CC=C(C=C1)C2=C(C=C(C=C2)C=C3C=CC=C3)C=C3C=CC=C3</chem>	<chem>CC1=CC=C(C=C1)C2=C(C=C(C=C2)C=C3C=CC=C3)C=C3C=CC=C3</chem>
52	<chem>CC1=CC=C(C=C1)C2=C(C=C(C=C2)Br)C=C3C=CC=C3</chem>	<chem>CC1=CC=C(C=C1)C2=C(C=C(C=C2)C=C3C=CC=C3)C=C3C=CC=C3</chem>	<chem>CC1=CC=C(C=C1)C2=C(C=C(C=C2)C=C3C=CC=C3)C=C3C=CC=C3</chem>
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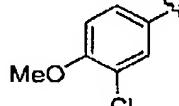
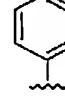
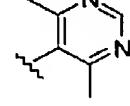
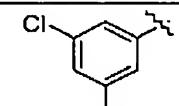
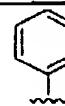
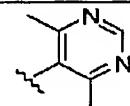
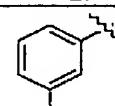
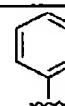
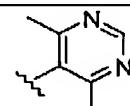
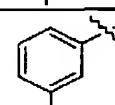
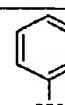
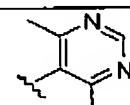
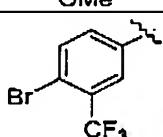
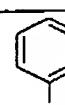
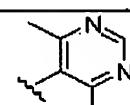
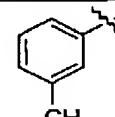
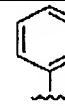
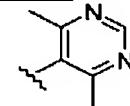
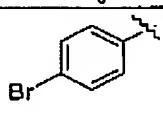
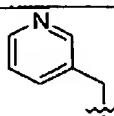
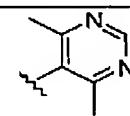
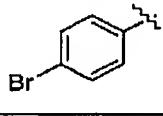
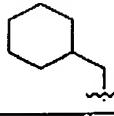
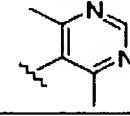
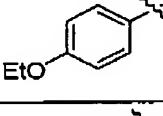
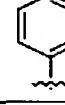
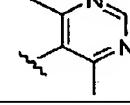
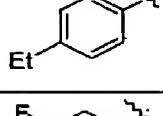
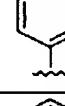
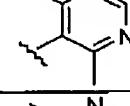
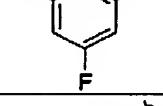
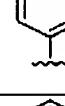
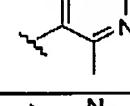
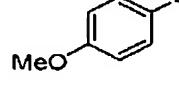
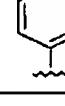
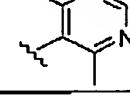
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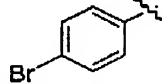
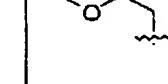
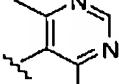
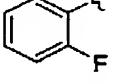
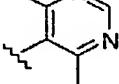
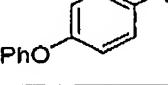
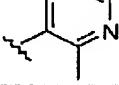
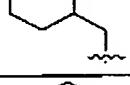
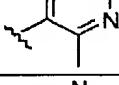
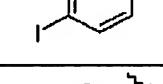
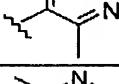
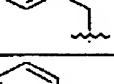
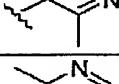
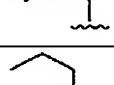
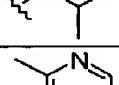
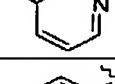
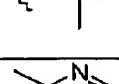
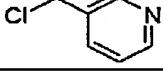
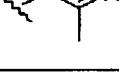
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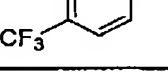
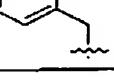
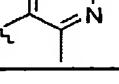
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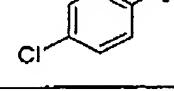
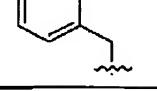
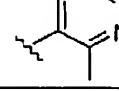
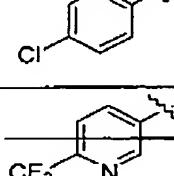
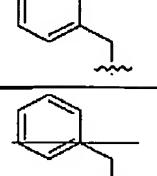
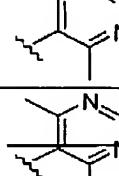
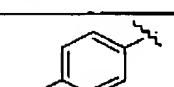
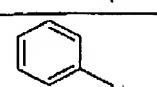
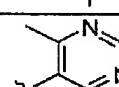
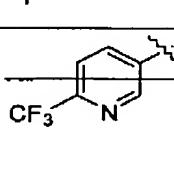
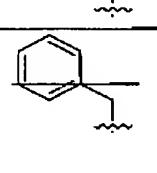
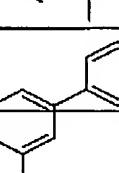
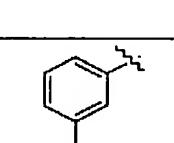
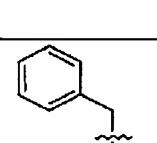
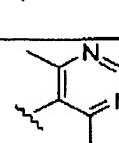
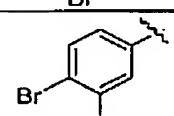
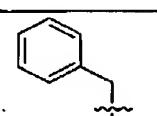
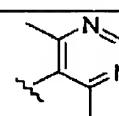
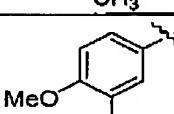
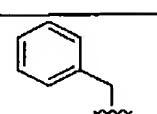
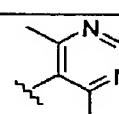
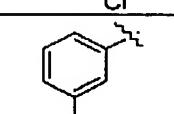
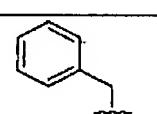
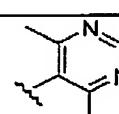
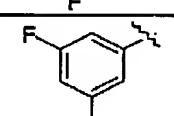
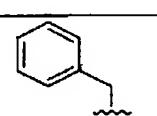
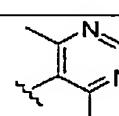
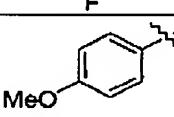
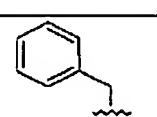
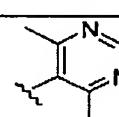
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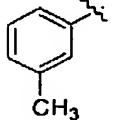
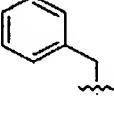
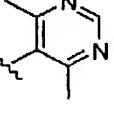
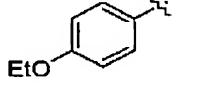
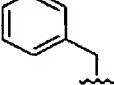
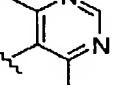
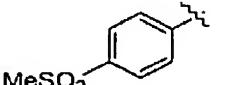
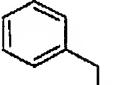
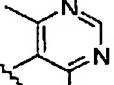
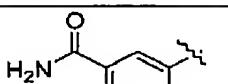
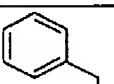
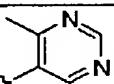
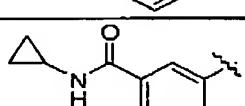
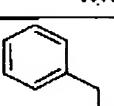
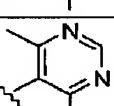
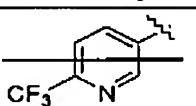
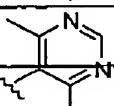
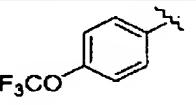
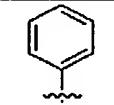
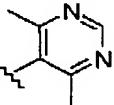
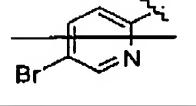
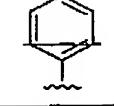
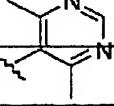
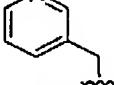
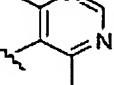
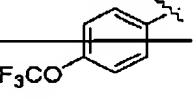
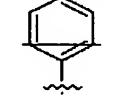
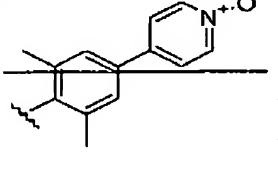
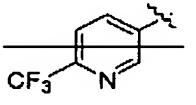
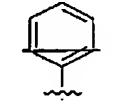
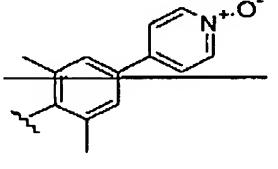
23. (Amended) A compound according to claim 22 wherein R<sup>1</sup>, R<sup>2</sup> and R<sup>3</sup> each represent:

#	R <sup>1</sup>	R <sup>2</sup>	R <sup>3</sup>
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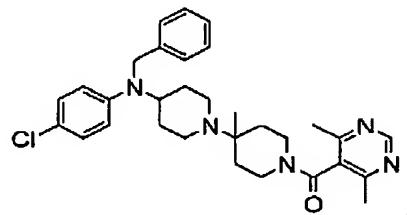
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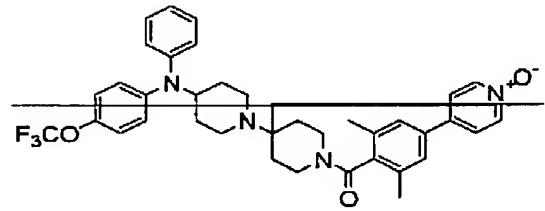
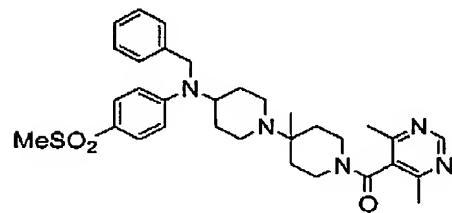
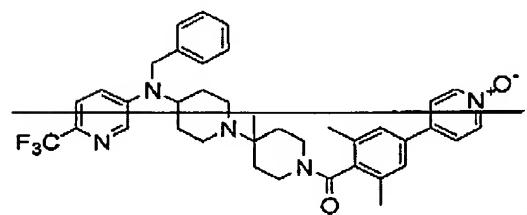
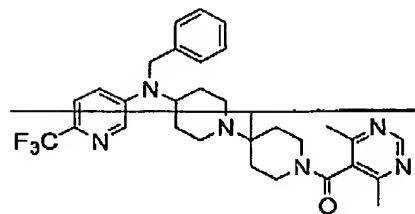
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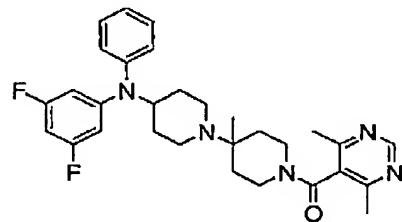
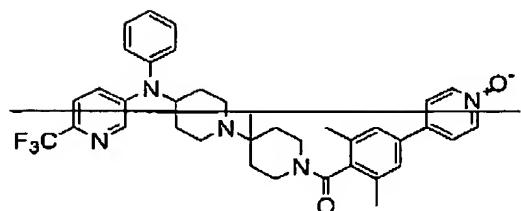
24. (Amended) A compound according to claim 23 represented by the structural formulae:



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25. (Previously presented) A pharmaceutical composition comprising one or more compounds of claim 21.

26. (Previously presented) A pharmaceutical composition comprising one or more compounds of claim 24.

27. (Previously presented) The pharmaceutical composition according to claim 25 further comprising one or more pharmaceutically acceptable carriers.

28. (Previously presented) The pharmaceutical composition according to claim 26 further comprising one or more pharmaceutically acceptable carriers.

29. (Previously presented) The pharmaceutical composition according to claim 25, wherein said pharmaceutical composition contains a therapeutically effective amount of said one or more compounds.

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30. (Previously presented) The pharmaceutical composition according to claim 26, wherein said pharmaceutical composition contains a therapeutically effective amount of said one or more compounds.

31-40. (Previously canceled)